Full wwPDB X-ray Structure Validation Report (i)

Nov 19, 2023 – 10:28 PM JST

PDB ID	:	7BXC
Title	:	Crystal structure of Ca_00311 form II
Authors	:	Fan, C.P.
Deposited on	:	2020-04-19
Resolution	:	2.25 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (i)) were used in the production of this report:

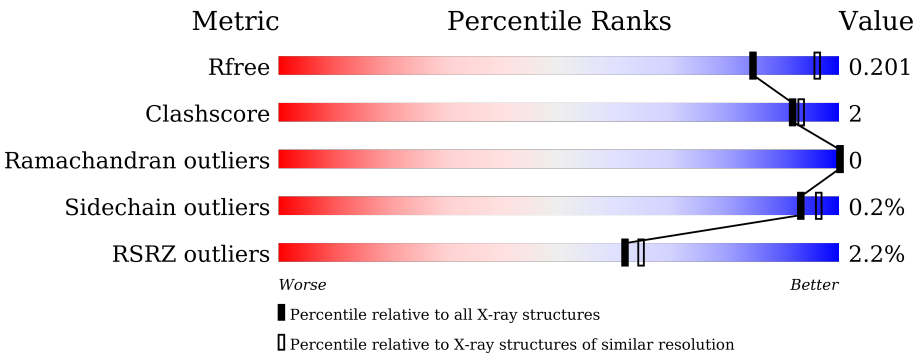
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	220	2% 98% •
1	B	220	2% 96% •
1	C	220	2% 93% 7%
1	D	220	95% 5%
1	E	220	% 95% 5%
1	F	220	2% 97% •

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Mol	Chain	Length	Quality of chain
1	G	220	% 97%
1	H	220	3% 97%
1	I	220	10% 94% 6%
1	J	220	3% 93% 5%
1	K	220	2% 96%
1	L	220	% 98%
1	M	220	% 98%
1	N	220	% 97%
1	O	220	2% 95% 5%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 28364 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

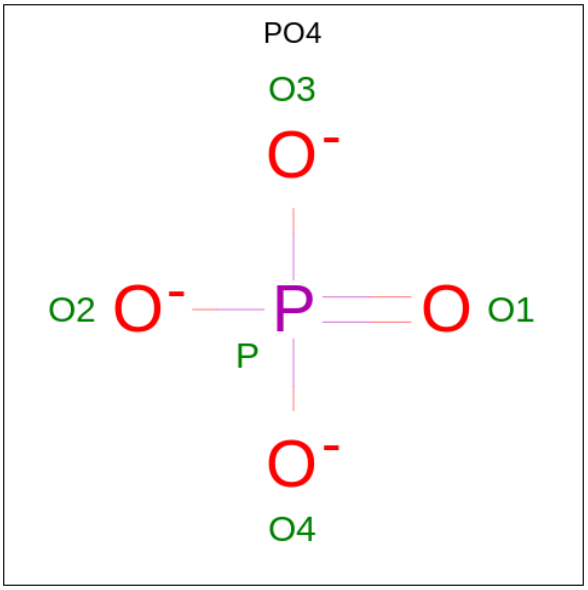
- Molecule 1 is a protein called GDSL-like Lipase/Acylhydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	220	1687	1060	293	317	17	0	0	0
1	B	220	1687	1060	293	317	17	0	0	0
1	C	220	1687	1060	293	317	17	0	0	0
1	D	220	1681	1057	290	317	17	0	0	0
1	E	220	1695	1064	293	321	17	0	0	0
1	F	220	1690	1062	293	318	17	0	0	0
1	G	220	1691	1062	293	319	17	0	0	0
1	H	220	1687	1060	293	317	17	0	0	0
1	I	220	1675	1054	292	312	17	0	0	0
1	J	218	1659	1045	287	310	17	0	0	0
1	K	220	1687	1060	293	317	17	0	0	0
1	L	220	1687	1060	292	318	17	0	0	0
1	M	220	1683	1058	293	315	17	0	0	0
1	N	220	1683	1058	293	315	17	0	0	0
1	O	220	1686	1060	293	316	17	0	0	0

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ALA	-	expression tag	UNP A0A381K2L8
B	0	ALA	-	expression tag	UNP A0A381K2L8
C	0	ALA	-	expression tag	UNP A0A381K2L8
D	0	ALA	-	expression tag	UNP A0A381K2L8
E	0	ALA	-	expression tag	UNP A0A381K2L8
F	0	ALA	-	expression tag	UNP A0A381K2L8
G	0	ALA	-	expression tag	UNP A0A381K2L8
H	0	ALA	-	expression tag	UNP A0A381K2L8
I	0	ALA	-	expression tag	UNP A0A381K2L8
J	0	ALA	-	expression tag	UNP A0A381K2L8
K	0	ALA	-	expression tag	UNP A0A381K2L8
L	0	ALA	-	expression tag	UNP A0A381K2L8
M	0	ALA	-	expression tag	UNP A0A381K2L8
N	0	ALA	-	expression tag	UNP A0A381K2L8
O	0	ALA	-	expression tag	UNP A0A381K2L8

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	P		
2	A	1	5	4	1	0	0
			Total	O	P		
2	A	1	5	4	1	0	0
			Total	O	P		
2	B	1	5	4	1	0	0
			Total	O	P		
2	B	1	5	4	1	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	P		
2	C	1	5	4	1	0	0
			Total	O	P		
2	C	1	5	4	1	0	0
			Total	O	P		
2	D	1	5	4	1	0	0
			Total	O	P		
2	D	1	5	4	1	0	0
			Total	O	P		
2	E	1	5	4	1	0	0
			Total	O	P		
2	E	1	5	4	1	0	0
			Total	O	P		
2	F	1	5	4	1	0	0
			Total	O	P		
2	F	1	5	4	1	0	0
			Total	O	P		
2	G	1	5	4	1	0	0
			Total	O	P		
2	G	1	5	4	1	0	0
			Total	O	P		
2	H	1	5	4	1	0	0
			Total	O	P		
2	H	1	5	4	1	0	0
			Total	O	P		
2	H	1	5	4	1	0	0
			Total	O	P		
2	I	1	5	4	1	0	0
			Total	O	P		
2	I	1	5	4	1	0	0
			Total	O	P		
2	J	1	5	4	1	0	0
			Total	O	P		
2	J	1	5	4	1	0	0
			Total	O	P		
2	K	1	5	4	1	0	0
			Total	O	P		
2	L	1	5	4	1	0	0
			Total	O	P		
2	L	1	5	4	1	0	0
			Total	O	P		
2	M	1	5	4	1	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	P		
2	M	1	5	4	1	0	0
2	N	1	5	4	1	0	0
2	N	1	5	4	1	0	0
2	O	1	5	4	1	0	0
2	O	1	5	4	1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	233	233	233	0	0
3	B	230	230	230	0	0
3	C	197	197	197	0	0
3	D	211	211	211	0	0
3	E	221	221	221	0	0
3	F	228	228	228	0	0
3	G	195	195	195	0	0
3	H	177	177	177	0	0
3	I	136	136	136	0	0
3	J	155	155	155	0	0
3	K	192	192	192	0	0
3	L	199	199	199	0	0
3	M	183	183	183	0	0
3	N	202	202	202	0	0

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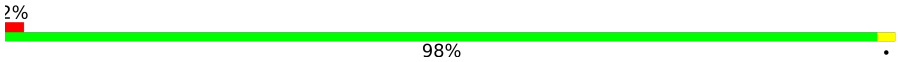
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	O	190	190	190	0	0

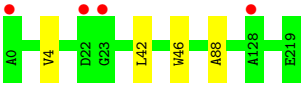
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: GDSL-like Lipase/Acylhydrolase

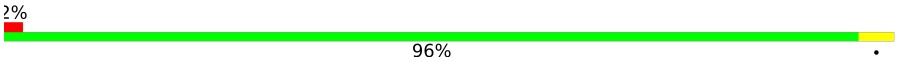
Chain A:





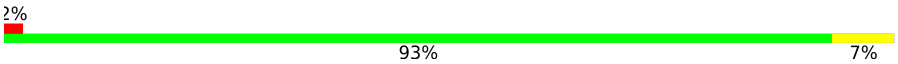
• Molecule 1: GDSL-like Lipase/Acylhydrolase

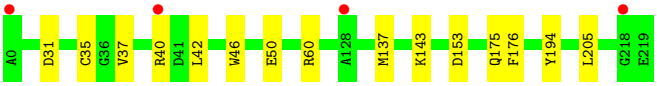
Chain B:



• Molecule 1: GDSL-like Lipase/Acylhydrolase

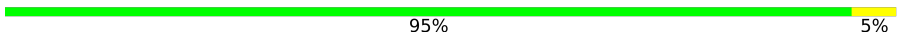
Chain C:





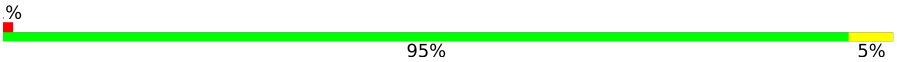
• Molecule 1: GDSL-like Lipase/Acylhydrolase

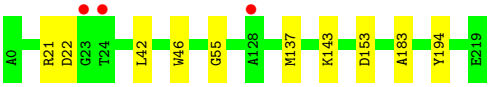
Chain D:



• Molecule 1: GDSL-like Lipase/Acylhydrolase

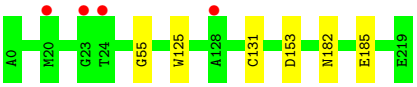
Chain E:





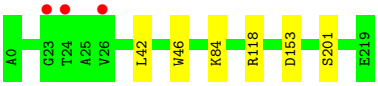
• Molecule 1: GDSL-like Lipase/Acylhydrolase

Chain F: 2% 97%



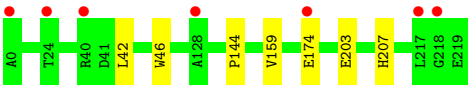
• Molecule 1: GDSL-like Lipase/Acylhydrolase

Chain G: % 97%



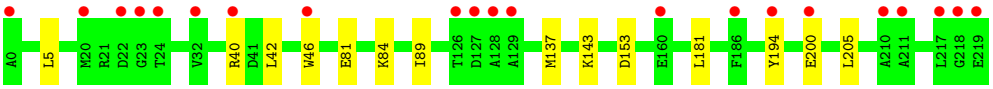
• Molecule 1: GDSL-like Lipase/Acylhydrolase

Chain H: 3% 97%



• Molecule 1: GDSL-like Lipase/Acylhydrolase

Chain I: 10% 94% 6%

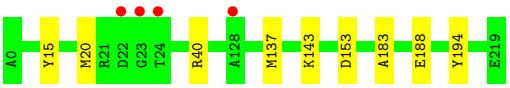


• Molecule 1: GDSL-like Lipase/Acylhydrolase

Chain J: 3% 93% 5%

• Molecule 1: GDSL-like Lipase/Acylhydrolase

Chain K: 2% 96%

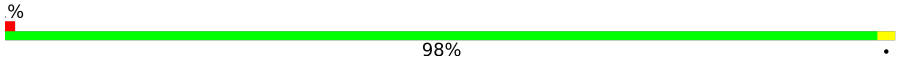


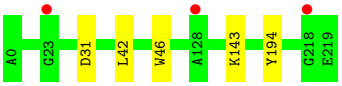
• Molecule 1: GDSL-like Lipase/Acylhydrolase

Chain L: % 98%

• Molecule 1: GDSL-like Lipase/Acylhydrolase

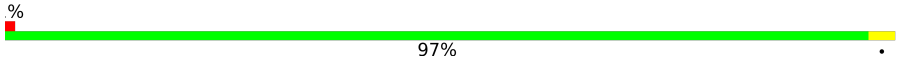
Chain M:

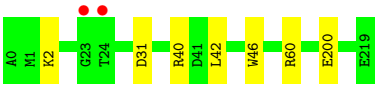




• Molecule 1: GDSL-like Lipase/Acylhydrolase

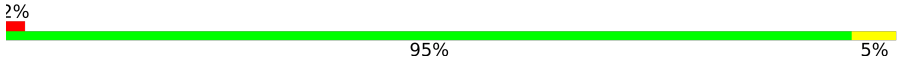
Chain N:





• Molecule 1: GDSL-like Lipase/Acylhydrolase

Chain O:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	326.44Å 109.96Å 152.98Å 90.00° 108.77° 90.00°	Depositor
Resolution (Å)	48.28 – 2.25 48.28 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.6 (48.28-2.25) 99.6 (48.28-2.25)	Depositor EDS
R_{merge}	0.24	Depositor
R_{sym}	0.24	Depositor
$< I/\sigma(I) >$ [1]	1.63 (at 2.24Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.165 , 0.201 0.165 , 0.201	Depositor DCC
R_{free} test set	12139 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	21.9	Xtriage
Anisotropy	0.461	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 54.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	28364	wwPDB-VP
Average B, all atoms (Å^2)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.82% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: PO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\|>5$	RMSZ	$\#\|Z\|>5$
1	A	0.27	0/1726	0.46	0/2342
1	B	0.26	0/1726	0.45	0/2342
1	C	0.26	0/1726	0.43	0/2342
1	D	0.26	0/1720	0.44	0/2335
1	E	0.26	0/1734	0.44	0/2352
1	F	0.26	0/1729	0.44	0/2347
1	G	0.26	0/1730	0.44	0/2347
1	H	0.25	0/1726	0.43	0/2342
1	I	0.25	0/1714	0.43	0/2327
1	J	0.26	0/1698	0.45	1/2309 (0.0%)
1	K	0.26	0/1726	0.44	0/2342
1	L	0.26	0/1726	0.43	0/2342
1	M	0.25	0/1722	0.43	0/2337
1	N	0.26	0/1722	0.43	0/2337
1	O	0.26	0/1725	0.44	0/2342
All	All	0.26	0/25850	0.44	1/35085 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	J	4	VAL	CG1-CB-CG2	6.61	121.47	110.90

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1687	0	1612	2	0
1	B	1687	0	1612	7	0
1	C	1687	0	1612	10	0
1	D	1681	0	1601	5	0
1	E	1695	0	1620	6	0
1	F	1690	0	1616	4	0
1	G	1691	0	1616	4	0
1	H	1687	0	1612	3	0
1	I	1675	0	1598	9	0
1	J	1659	0	1577	7	0
1	K	1687	0	1612	6	0
1	L	1687	0	1610	3	0
1	M	1683	0	1608	2	0
1	N	1683	0	1608	6	0
1	O	1686	0	1612	6	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
2	C	10	0	0	0	0
2	D	10	0	0	0	0
2	E	10	0	0	1	0
2	F	10	0	0	1	0
2	G	10	0	0	0	0
2	H	15	0	0	0	0
2	I	10	0	0	0	0
2	J	10	0	0	1	0
2	K	5	0	0	0	0
2	L	10	0	0	0	0
2	M	10	0	0	0	0
2	N	10	0	0	0	0
2	O	10	0	0	0	0
3	A	233	0	0	0	0
3	B	230	0	0	2	0
3	C	197	0	0	0	0
3	D	211	0	0	0	0
3	E	221	0	0	0	0
3	F	228	0	0	0	0
3	G	195	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	H	177	0	0	0	0
3	I	136	0	0	0	0
3	J	155	0	0	0	0
3	K	192	0	0	2	0
3	L	199	0	0	0	0
3	M	183	0	0	0	0
3	N	202	0	0	1	0
3	O	190	0	0	0	1
All	All	28364	0	24126	80	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 2.

All (80) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:188:GLU:OE1	3:K:401:HOH:O	1.81	0.95
1:I:137:MET:HE1	1:I:205:LEU:HD11	1.69	0.74
1:O:31:ASP:HB3	1:O:40:ARG:HH21	1.57	0.68
1:C:137:MET:HE1	1:C:205:LEU:HD11	1.74	0.67
1:F:55:GLY:N	2:F:302:PO4:O2	2.30	0.65
1:I:40:ARG:HG2	1:I:40:ARG:HH11	1.62	0.64
1:O:137:MET:HE3	1:O:205:LEU:HD11	1.81	0.61
1:E:55:GLY:N	2:E:301:PO4:O2	2.36	0.57
1:J:4:VAL:HG23	1:J:88:ALA:HB3	1.86	0.56
1:J:171:HIS:O	1:J:175:GLN:HG2	2.07	0.55
1:N:60:ARG:NH2	3:N:401:HOH:O	2.33	0.54
1:O:15:TYR:CZ	1:O:20:MET:HG3	2.43	0.54
1:G:118:ARG:NH2	3:G:402:HOH:O	2.41	0.53
1:J:83:HIS:ND1	2:J:302:PO4:O4	2.43	0.52
1:K:137:MET:HE1	1:K:183:ALA:HB2	1.92	0.51
1:G:153:ASP:OD1	1:G:153:ASP:N	2.39	0.51
1:B:26:VAL:HG22	3:B:528:HOH:O	2.09	0.51
1:K:40:ARG:NH1	3:K:404:HOH:O	2.44	0.50
1:C:175:GLN:HG3	1:C:176:PHE:HD1	1.76	0.50
1:C:37:VAL:HG22	1:C:40:ARG:HH21	1.76	0.49
1:F:153:ASP:OD1	1:F:153:ASP:N	2.39	0.49
1:K:143:LYS:HB2	1:K:194:TYR:HB3	1.94	0.49
1:C:175:GLN:HG3	1:C:176:PHE:CD1	2.48	0.49
1:C:31:ASP:O	1:C:40:ARG:NH2	2.46	0.48
1:N:200:GLU:H	1:N:200:GLU:CD	2.16	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:42:LEU:HB3	1:E:46:TRP:CG	2.49	0.48
1:M:42:LEU:HB3	1:M:46:TRP:CG	2.49	0.47
1:H:42:LEU:HB3	1:H:46:TRP:CG	2.50	0.47
1:D:175:GLN:HG3	1:D:176:PHE:CD1	2.50	0.47
1:J:42:LEU:HB3	1:J:46:TRP:CG	2.50	0.47
1:E:21:ARG:HD2	1:E:22:ASP:OD1	2.15	0.46
1:A:4:VAL:HG22	1:A:88:ALA:HB3	1.97	0.46
1:E:143:LYS:HB2	1:E:194:TYR:HB3	1.98	0.46
1:F:125:TRP:CD1	1:F:131:CYS:HB2	2.51	0.46
1:N:2:LYS:HD3	1:N:46:TRP:CE2	2.51	0.46
1:D:143:LYS:HB2	1:D:194:TYR:HB3	1.98	0.46
1:M:143:LYS:HB2	1:M:194:TYR:HB3	1.98	0.46
1:D:42:LEU:HB3	1:D:46:TRP:CG	2.51	0.46
1:N:42:LEU:HB3	1:N:46:TRP:CG	2.51	0.46
1:C:35:CYS:HB2	1:C:50:GLU:OE2	2.15	0.45
1:O:137:MET:HE1	1:O:183:ALA:HB2	1.97	0.45
1:D:21:ARG:HD2	1:D:22:ASP:OD1	2.17	0.45
1:J:15:TYR:CZ	1:J:20:MET:HG3	2.52	0.45
1:D:171:HIS:O	1:D:174:GLU:HG2	2.17	0.45
1:B:42:LEU:HB3	1:B:46:TRP:CG	2.52	0.45
1:I:143:LYS:HB2	1:I:194:TYR:HB3	1.97	0.45
1:O:134:ILE:HB	1:O:178:CYS:HB2	1.98	0.45
1:A:42:LEU:HB3	1:A:46:TRP:CG	2.51	0.45
1:B:118:ARG:NH2	3:B:402:HOH:O	2.40	0.44
1:C:37:VAL:HG22	1:C:40:ARG:NH2	2.33	0.44
1:G:42:LEU:HB3	1:G:46:TRP:CG	2.53	0.44
1:H:203:GLU:OE2	1:H:207:HIS:NE2	2.50	0.44
1:I:200:GLU:H	1:I:200:GLU:CD	2.21	0.44
1:K:153:ASP:OD1	1:K:153:ASP:N	2.41	0.44
1:E:153:ASP:OD1	1:E:153:ASP:N	2.43	0.43
1:G:84:LYS:NZ	3:G:404:HOH:O	2.49	0.43
1:K:15:TYR:CZ	1:K:20:MET:HG3	2.53	0.43
1:C:153:ASP:OD1	1:C:153:ASP:N	2.39	0.43
1:I:153:ASP:OD1	1:I:153:ASP:N	2.40	0.43
1:I:42:LEU:HB3	1:I:46:TRP:CG	2.54	0.43
1:J:41:ASP:OD2	1:J:207:HIS:ND1	2.35	0.43
1:B:175:GLN:HG3	1:B:176:PHE:CD1	2.54	0.42
1:B:175:GLN:HG3	1:B:176:PHE:HD1	1.84	0.42
1:I:81:GLU:OE2	1:I:84:LYS:HE2	2.20	0.42
1:L:153:ASP:OD1	1:L:153:ASP:N	2.44	0.42
1:C:42:LEU:HB3	1:C:46:TRP:CG	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:5:LEU:HD23	1:I:89:ILE:HD12	2.02	0.42
1:N:31:ASP:N	1:N:31:ASP:OD1	2.52	0.42
1:N:31:ASP:HB2	1:N:40:ARG:HH22	1.85	0.42
1:O:125:TRP:CD1	1:O:131:CYS:HB2	2.54	0.42
1:L:143:LYS:HB2	1:L:194:TYR:HB3	2.02	0.41
1:C:143:LYS:HB2	1:C:194:TYR:HB3	2.03	0.41
1:F:182:ASN:HB3	1:F:185:GLU:HG3	2.03	0.41
1:B:137:MET:HA	1:B:181:LEU:O	2.20	0.41
1:H:144:PRO:HA	1:H:159:VAL:HG21	2.02	0.41
1:E:137:MET:HE1	1:E:183:ALA:HB2	2.03	0.40
1:L:144:PRO:HA	1:L:159:VAL:HG21	2.02	0.40
1:B:153:ASP:OD1	1:B:153:ASP:N	2.43	0.40
1:J:143:LYS:HB2	1:J:194:TYR:HB3	2.03	0.40
1:I:137:MET:HA	1:I:181:LEU:O	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:O:408:HOH:O	3:O:565:HOH:O[2_556]	2.16	0.04

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	218/220 (99%)	209 (96%)	9 (4%)	0	100	100
1	B	218/220 (99%)	209 (96%)	9 (4%)	0	100	100
1	C	218/220 (99%)	209 (96%)	9 (4%)	0	100	100
1	D	218/220 (99%)	210 (96%)	8 (4%)	0	100	100
1	E	218/220 (99%)	210 (96%)	8 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	218/220 (99%)	209 (96%)	9 (4%)	0	100	100
1	G	218/220 (99%)	209 (96%)	9 (4%)	0	100	100
1	H	218/220 (99%)	209 (96%)	9 (4%)	0	100	100
1	I	218/220 (99%)	209 (96%)	9 (4%)	0	100	100
1	J	216/220 (98%)	207 (96%)	9 (4%)	0	100	100
1	K	218/220 (99%)	210 (96%)	8 (4%)	0	100	100
1	L	218/220 (99%)	209 (96%)	9 (4%)	0	100	100
1	M	218/220 (99%)	209 (96%)	9 (4%)	0	100	100
1	N	218/220 (99%)	209 (96%)	9 (4%)	0	100	100
1	O	218/220 (99%)	209 (96%)	9 (4%)	0	100	100
All	All	3268/3300 (99%)	3136 (96%)	132 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	172/176 (98%)	172 (100%)	0	100	100
1	B	172/176 (98%)	172 (100%)	0	100	100
1	C	172/176 (98%)	171 (99%)	1 (1%)	86	91
1	D	171/176 (97%)	171 (100%)	0	100	100
1	E	174/176 (99%)	174 (100%)	0	100	100
1	F	173/176 (98%)	173 (100%)	0	100	100
1	G	173/176 (98%)	172 (99%)	1 (1%)	86	91
1	H	172/176 (98%)	171 (99%)	1 (1%)	86	91
1	I	169/176 (96%)	169 (100%)	0	100	100
1	J	168/176 (96%)	168 (100%)	0	100	100
1	K	172/176 (98%)	172 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	172/176 (98%)	172 (100%)	0	100	100
1	M	171/176 (97%)	170 (99%)	1 (1%)	86	91
1	N	171/176 (97%)	171 (100%)	0	100	100
1	O	172/176 (98%)	172 (100%)	0	100	100
All	All	2574/2640 (98%)	2570 (100%)	4 (0%)	93	96

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	60	ARG
1	G	201	SER
1	H	174	GLU
1	M	31	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

30 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	PO4	O	302	-	4,4,4	0.96	0	6,6,6	0.46	0
2	PO4	F	302	-	4,4,4	1.05	0	6,6,6	0.35	0
2	PO4	N	301	-	4,4,4	0.95	0	6,6,6	0.44	0
2	PO4	E	301	-	4,4,4	1.05	0	6,6,6	0.39	0
2	PO4	L	301	-	4,4,4	0.91	0	6,6,6	0.48	0
2	PO4	C	302	-	4,4,4	0.93	0	6,6,6	0.33	0
2	PO4	K	301	-	4,4,4	0.88	0	6,6,6	0.51	0
2	PO4	A	302	-	4,4,4	0.97	0	6,6,6	0.46	0
2	PO4	L	302	-	4,4,4	0.94	0	6,6,6	0.51	0
2	PO4	O	301	-	4,4,4	0.90	0	6,6,6	0.51	0
2	PO4	I	301	-	4,4,4	0.90	0	6,6,6	0.44	0
2	PO4	F	301	-	4,4,4	0.98	0	6,6,6	0.45	0
2	PO4	C	301	-	4,4,4	0.97	0	6,6,6	0.38	0
2	PO4	G	301	-	4,4,4	0.92	0	6,6,6	0.37	0
2	PO4	B	302	-	4,4,4	0.94	0	6,6,6	0.30	0
2	PO4	J	301	-	4,4,4	0.91	0	6,6,6	0.39	0
2	PO4	H	301	-	4,4,4	0.98	0	6,6,6	0.45	0
2	PO4	A	301	-	4,4,4	1.01	0	6,6,6	0.45	0
2	PO4	G	302	-	4,4,4	0.95	0	6,6,6	0.48	0
2	PO4	E	302	-	4,4,4	0.99	0	6,6,6	0.39	0
2	PO4	M	302	-	4,4,4	1.01	0	6,6,6	0.55	0
2	PO4	N	302	-	4,4,4	0.92	0	6,6,6	0.39	0
2	PO4	I	302	-	4,4,4	0.92	0	6,6,6	0.52	0
2	PO4	J	302	-	4,4,4	0.97	0	6,6,6	0.43	0
2	PO4	B	301	-	4,4,4	0.97	0	6,6,6	0.43	0
2	PO4	H	302	-	4,4,4	0.93	0	6,6,6	0.49	0
2	PO4	H	303	-	4,4,4	0.95	0	6,6,6	0.38	0
2	PO4	D	301	-	4,4,4	0.96	0	6,6,6	0.41	0
2	PO4	M	301	-	4,4,4	0.97	0	6,6,6	0.44	0
2	PO4	D	302	-	4,4,4	0.97	0	6,6,6	0.45	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	302	PO4	1	0
2	E	301	PO4	1	0
2	J	302	PO4	1	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	220/220 (100%)	-0.33	4 (1%)	68	71	9, 19, 37, 53	0
1	B	220/220 (100%)	-0.33	4 (1%)	68	71	11, 20, 37, 59	0
1	C	220/220 (100%)	-0.01	4 (1%)	68	71	12, 24, 46, 58	0
1	D	220/220 (100%)	-0.32	0	100	100	14, 21, 38, 65	0
1	E	220/220 (100%)	-0.24	3 (1%)	75	77	12, 20, 40, 64	0
1	F	220/220 (100%)	-0.12	4 (1%)	68	71	11, 21, 41, 52	0
1	G	220/220 (100%)	-0.16	3 (1%)	75	77	14, 24, 45, 64	0
1	H	220/220 (100%)	0.01	7 (3%)	47	50	15, 28, 51, 64	0
1	I	220/220 (100%)	0.50	21 (9%)	8	8	20, 36, 60, 79	0
1	J	218/220 (99%)	0.10	6 (2%)	53	55	21, 34, 53, 66	0
1	K	220/220 (100%)	-0.20	4 (1%)	68	71	15, 23, 45, 59	0
1	L	220/220 (100%)	-0.28	2 (0%)	84	85	16, 25, 43, 61	0
1	M	220/220 (100%)	-0.18	3 (1%)	75	77	17, 27, 45, 57	0
1	N	220/220 (100%)	-0.29	2 (0%)	84	85	16, 23, 40, 56	0
1	O	220/220 (100%)	-0.22	5 (2%)	60	63	17, 26, 45, 67	0
All	All	3298/3300 (99%)	-0.14	72 (2%)	62	65	9, 24, 46, 79	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	23	GLY	5.2
1	E	128	ALA	4.9
1	O	128	ALA	4.8
1	I	0	ALA	4.0
1	H	128	ALA	3.7
1	N	24	THR	3.6
1	B	24	THR	3.6

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Mol	Chain	Res	Type	RSRZ
1	H	0	ALA	3.5
1	H	218	GLY	3.5
1	O	23	GLY	3.5
1	I	24	THR	3.4
1	C	218	GLY	3.4
1	K	23	GLY	3.2
1	I	23	GLY	3.2
1	I	217	LEU	3.2
1	N	23	GLY	3.2
1	G	23	GLY	3.1
1	I	218	GLY	3.1
1	I	32	VAL	3.0
1	I	22	ASP	2.9
1	G	24	THR	2.9
1	I	128	ALA	2.9
1	C	40	ARG	2.9
1	F	24	THR	2.9
1	E	24	THR	2.9
1	H	217	LEU	2.9
1	A	0	ALA	2.8
1	I	46	TRP	2.8
1	B	23	GLY	2.7
1	M	218	GLY	2.7
1	J	24	THR	2.7
1	B	26	VAL	2.7
1	O	22	ASP	2.7
1	K	22	ASP	2.7
1	J	23	GLY	2.6
1	J	66	LEU	2.6
1	J	144	PRO	2.6
1	K	24	THR	2.6
1	M	128	ALA	2.6
1	I	40	ARG	2.5
1	I	194	TYR	2.5
1	I	219	GLU	2.5
1	I	20	MET	2.5
1	A	22	ASP	2.4
1	M	23	GLY	2.4
1	I	186	PHE	2.4
1	A	23	GLY	2.3
1	O	24	THR	2.3
1	I	211	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	H	40	ARG	2.3
1	I	129	ALA	2.3
1	B	22	ASP	2.3
1	I	127	ASP	2.3
1	H	24	THR	2.3
1	C	128	ALA	2.2
1	I	210	ALA	2.2
1	J	0	ALA	2.2
1	K	128	ALA	2.2
1	I	126	THR	2.2
1	H	174	GLU	2.2
1	L	128	ALA	2.1
1	I	200	GLU	2.1
1	E	23	GLY	2.1
1	L	19	GLY	2.1
1	A	128	ALA	2.1
1	F	20	MET	2.1
1	F	128	ALA	2.1
1	G	26	VAL	2.1
1	O	20	MET	2.0
1	I	160	GLU	2.0
1	C	0	ALA	2.0
1	J	21	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	PO4	F	302	5/5	0.95	0.19	23,24,42,45	0
2	PO4	E	301	5/5	0.96	0.18	27,28,34,39	0
2	PO4	D	301	5/5	0.96	0.16	23,28,34,42	0
2	PO4	I	301	5/5	0.96	0.14	45,49,52,59	0
2	PO4	J	302	5/5	0.96	0.21	46,48,56,58	0
2	PO4	C	301	5/5	0.97	0.24	31,36,45,47	0
2	PO4	H	303	5/5	0.97	0.21	32,41,44,45	0
2	PO4	H	301	5/5	0.98	0.14	31,34,41,42	0
2	PO4	F	301	5/5	0.98	0.23	22,37,42,47	0
2	PO4	M	302	5/5	0.98	0.22	32,39,43,48	0
2	PO4	N	301	5/5	0.98	0.13	35,36,40,40	0
2	PO4	G	301	5/5	0.99	0.09	19,21,26,26	0
2	PO4	G	302	5/5	0.99	0.14	28,29,39,43	0
2	PO4	B	301	5/5	0.99	0.11	27,29,39,42	0
2	PO4	H	302	5/5	0.99	0.08	22,25,25,27	0
2	PO4	D	302	5/5	0.99	0.12	26,32,36,37	0
2	PO4	B	302	5/5	0.99	0.09	12,17,19,25	0
2	PO4	I	302	5/5	0.99	0.10	24,28,31,35	0
2	PO4	J	301	5/5	0.99	0.09	29,30,34,38	0
2	PO4	E	302	5/5	0.99	0.14	20,30,33,33	0
2	PO4	K	301	5/5	0.99	0.08	22,24,25,27	0
2	PO4	L	301	5/5	0.99	0.08	20,21,22,23	0
2	PO4	L	302	5/5	0.99	0.13	30,30,37,52	0
2	PO4	M	301	5/5	0.99	0.10	15,19,23,30	0
2	PO4	A	302	5/5	0.99	0.14	22,24,36,37	0
2	PO4	C	302	5/5	0.99	0.09	15,17,19,23	0
2	PO4	N	302	5/5	0.99	0.07	19,20,23,26	0
2	PO4	O	301	5/5	0.99	0.08	23,23,27,27	0
2	PO4	O	302	5/5	0.99	0.15	37,37,45,47	0
2	PO4	A	301	5/5	1.00	0.10	13,15,19,19	0

6.5 Other polymers (i)

There are no such residues in this entry.